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AFWL-TR-66-108, Vol. III

AFWL-TR  
66-108  
Vol. III



# NUCLEAR EXPLOSION INTERACTION STUDIES

Volume III

Miscellaneous Code Development

K. D. Pyatt, Jr., et al.

General Atomic Division  
General Dynamics Corporation  
Special Nuclear Effects Laboratory  
San Diego, California  
Contract AF 29(601)-7035

TECHNICAL REPORT NO. AFWL-TR-66-108, Vol. III

May 1967

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FOREWORD

This report was prepared by General Atomic Division, General Dynamics Corporation, San Diego, California, under Contract AF29(601)-7035. The research was funded by DASA under Program Element 6.16.46.01D, Project 5710, Subtask 07.002, and by ARPA Order 313, Program Element 6.25.03.01.R.

Inclusive dates of research were 22 July 1965 to 21 July 1966. The report was submitted 28 April 1967 by the AFWL Project Officer, Maj George Spillman (WLRT). The contractor's report number is GA-7370.


This final report on Nuclear Explosion Interaction Studies is being published in four volumes. The volume titles are as follows: Volume I, Methods for Analysis of Radiative Transfer; Volume II, Methods for Analysis of Thermal Phenomena; Volume III, Miscellaneous Code Development; and Volume IV, Phenomenology Studies (classified SECRET/RESTRICTED DATA).

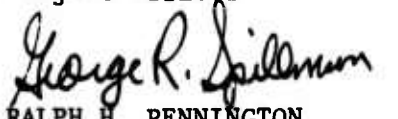
The first three volumes are devoted, respectively, to theoretical studies and computer code development in radiative transfer, thermal phenomena, and miscellaneous efforts related to various other aspects of the work. The fourth volume, which is classified, contains the results of applications of these techniques, and of those previously developed, to the study of fireball growth and the interaction of laser radiation with materials.

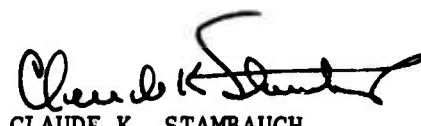
The NEIS program is long-range, and most of the projects described in this report are in an incomplete state of development. This is due in part to the nature of the existing computer programs themselves, which continue in a state of development as long as they are in use, and in part to the time scale involved in bringing new programs to a state of capability for solving real problems.

General Atomic staff personnel contributing to the research include J.H. Alexander, C.R. Dismukes, R. Durstenfeld, R.S. Engelmores, B.E. Freeman, W.B. Lindley, J.T. Palmer, K.D. Pyatt, L.L. Reed, L.M. Schalit, J.R. Triplett, and numerous others. The cooperation of Col R.H. Pennington, Maj G.R. Spillman, Lt B.S. Chambers, III, Lt N.D. Morgan, Lt R.A. Howerton, Dr. P.V. Avizonis, and Mr. D.W. Lane of the Air Force Weapons Laboratory is gratefully acknowledged.

This report has been reviewed and is approved.

  
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ABSTRACT

A group of SPUTTER routines has been modified or developed to execute in a more nearly accurate and efficient manner the task of "rezoning" the Lagrangian zone structure in a specified region of the system being analyzed. The automatic rezoning routine AUTORZ has been expanded to provide for maintaining a prescribed spatial definition within a region. The control routines REZONE, CONVER, and ZONE have been modified to make use of the new routines ZMS and ZRAD, which accomplish conversion between space and mass coordinates, and a special procedure PROFL is described which performs energy interpolation in certain cases. To facilitate maintenance of complex computer codes such as SPUTTER, OIL, and HECTIC, two FORTRAN analysis routines (FAR, Programs One and Two) have been developed and are nearly complete at this time. These routines are designed for replacement of FORTRAN symbol strings by others, and for compilation of complete cross reference listings of variable and subroutine references. Editing routines designed for specific use with the data-link terminal facilities now installed at the General Atomic Special Nuclear Effects Laboratory (SNEL) and at AFWL are also described.

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## SECTION I

### GENERATOR-REZONE ROUTINES FOR SPUTTER

In response to the needs of several applications, extensive modifications have been made in most of the SPUTTER routines to establish or revise the zone structure which partitions the material being studied into convenient slices for computational purposes. The more fundamental of these changes are discussed below.

#### 1.1. AUTORZ (Automatic Rezone Specification)

##### 1.1.1. Introduction

The purpose of this routine is to maintain, during the course of a calculation, an adequately fine-grained zone structure in those portions of the material being studied which are of principal physical interest, and at the same time avoid wasteful use of small zones in other portions. The routine is presently adapted to a restricted class of problems, namely, those in which an ablation front is moving into solid material on the left, with vapor expanding to the right. Only a few large zones are needed to describe the conduction front in the solid material, whereas the zone at the ablation front and the zones in the vapor for a certain distance outward must be much smaller if the objectives of the calculation are to be achieved. The outermost portions of the vapor are of less interest and can be described by a less detailed zone structure.

Since SPUTTER is a Lagrangian code, the zone boundaries move with the material and the problem is particularly serious. Even in an Eulerian code, however, a similar problem arises unless a moving coordinate system can be used with respect to which the ablation front is stationary. The

techniques described in this section may easily be generalized to accommodate problems associated with a wide variety of situations other than those associated with ablation; near shock and radiation fronts, for example, zone definition may be required which is difficult or impossible to maintain without repeated redefinition of the zone structure.

AUTORZ is called at the beginning of each cycle if the flag EKLM (19) is nonzero. The AUTORZ option control, THETAK (58), must also be nonzero or there is an immediate return.

#### 1.1.2. The Solid Rezone Procedure

In the solid material it is necessary to set up a few zones with a very large mass available for ablation and for containing the conduction front, and to split these, one at a time, into small zones as the ablation front advances into the solid. It is assumed that the large zones initially generated were equal in mass, although this is not strictly essential. The splitting ratio is an input quantity, THETAK(53), and is typically set at a number like 10.0 or 20.0. It must be an integer but has REAL type. The index of the boundary between the large zones and the small ones formed by the splitting operation is stored as THETAK(54); this is reduced by unity each time a large zone is split. It is usually best to generate the problem with a "lambda region" of small zones already present, in which case THETAK(54) should be initially set to the left boundary index of this lambda region.

The solid rezone is executed when IA, the ablation front index, has been reduced to the value THETAK(54)+THETAK(57). THETAK(57) is an input specification, usually of the order of 3 to 10, which establishes a minimum for the number of small zones behind the ablation front. This is advisable for reasons involving the conduction and source deposition calculations in some applications.

#### 1.1.3. The Inner Vapor Rezone Procedure

In the analysis of certain experimental arrangements, such as plasma probe measurements, good zone definition must be maintained within a

region of space defined in the Eulerian sense, i.e., in terms of fixed coordinates rather than moving Lagrangian zone boundaries. The region of interest is assumed to extend from the ablation front on the left to a coordinate position specified by the input in THETAK(59). Within this region, the desired zone width is specified by THETAK(60). If THETAK(59)=0 or if the leading edge of the vapor has not yet passed beyond THETAK(59), no attempt to rezone the inner vapor is made. The index K of the zone which contains the outer boundary THETAK(59) of the region of interest and the index J of the outermost zone within this region whose width is less than THETAK(60) are then identified. Unless the solid rezone is being done at the same time, the inner vapor rezone is omitted when the width of zone K is less than 1.5 times THETAK(60). If the solid rezone is to be done anyway, the inner vapor rezone is also done provided that there is at least one zone between J and K.

The inner vapor rezone is specified as follows: left boundary index  $IL=J+1$ , right boundary index  $IR=K+1$ , left zone width  $VAL(2)=THETAK(60)$ , number of zones  $VAL(5)=$ one more than the minimum number for which none of the new zones would exceed THETAK(60) in width. This specification generates new zones whose width decreases slightly toward the right, a desirable situation since it compensates in advance for later expansion of the zones.

#### 1.1.4. The Outer Vapor Rezone

In order to prevent the total number of zones from increasing too rapidly due to repeated solid rezoning, the outer part of the vapor is rezoned periodically to increase zone size in this region. The outer region extends from the outer boundary of the "inner vapor region" to the leading edge of the vapor, R(IB). The inner vapor region consists of the rezone region defined in the preceding paragraph, if such a rezone is being done; otherwise it consists of a definite (input) number THETAK(55) of zones counted out from the ablation front. The outer vapor rezone is normally done whenever the solid or inner vapor rezones are done; however, if the inner vapor is

being rezoned, the outer region must have a minimum of THETAK(55) zones in it or it is not considered worthwhile.

The outer region is rezoned as follows: left-most zone mass VAL(2) = the mass of the outermost of the inner zones, ratio of each zone mass to its neighbor on the left VAL(4)=THETAK(56). (In order to reduce the number of zones, THETAK(56) must be greater than 1.)

#### 1.1.5. Additional Controls and Flags

If THETAK(58)=1., only solid rezones can be done; if THETAK(58)=-1., solid rezones are possible only on cycles for which both inner and outer vapor rezones are also done. For THETAK(58)=2., there are no controls beyond those listed in the preceding paragraphs. If a solid rezone is to be done, the temperature profile flag is set (this is a minus sign on the right boundary index IR for the solid rezone specification). If any rezones are to be done, the linkage flag S16 is set to -1, indicating to the rezone programs that they are to rezone from data in RK, and MAIN1 is then called.

#### 1.1.6. The RK Array

The RK array contains specifications for a maximum of nine rezone regions, which the rezone program will process in the order of placement in the RK array. Thus, AUTORIZ specifies that the solid rezone is to be done after the vapor rezones by listing its specifications after those for the vapor rezones (if any).

RK(1) contains the number of rezone regions specified in the array. Each specification consists of 11 words, i.e., RK(2)-RK(12), RK(13)-RK(23), etc. These words are, respectively:

IL	Left boundary index of region (REAL type)
IR	Right boundary index of region (REAL type)
VAL(2)	Mass (if positive) or width (if negative) of left-most zone
VAL(3)	Mass (if positive) or width (if negative) of right-most zone

VAL(4)	Width ratio (if VAL(6)=1) or mass ratio (if VAL(6)=2) of each zone to its neighbor on the left
VAL(5)	Number of zones to be generated in the region
VAL(6)	Mass-space switch (see VAL(4) )
VAL(7)	Material identification number
VAL(8)	Multiplicative factor for opacity
VAL(9)	$\bar{A}$ (mean atomic mass number)
VAL(10)	ZMEAN (mean atomic number)

Of these 11 parameters, those required are the first two, at least two of the next four, and the seventh (VAL(6)). The others can either be calculated from those given or retain their previous values unchanged. The AUTORZ routine always leaves the last four words unspecified.

## 1.2. REZONE (REZONE CONTROL ROUTINE)

On entry to this routine, rezone specifications have been placed in the RK array either by AUTORZ or by the ZONERD routine which reads the specifications from card input. The number of such specifications is given by RK(1), and the entire procedure described below is repeated for each specification in turn. The format for the specifications is given in Section 1.1.6.

The rezone procedure will modify information within the rezone region defined by the boundary index range  $IL \leq I \leq IR$ . Unmodified values must first be saved, as follows:



		<u>Quantity</u>	<u>Save location</u>	
For all I, $IL \leq I \leq IR$		Velocities, RD(I)	RDTAB(N)	in /NAMEC/ COMMON
		Densities, 1./SV(I)	RHOTAB(N)	
		Specific energies, E(I)	THETAB(N)	
		Coordinates, R(I)	RTABRD(N)	
		Coordinates, R(I)	RTABRO(N)	
		Coordinates, R(I)	RTABTH(N)	
		Mean ionic charges, FEW(I)	CAR(N)	in blank COMMON
		$Z^1$ , X6(I)	CAPAC(N)	
		$Z^2$ , BC(I)	CHIC(N)	
		$Z^3$ , BR(I)	CHIR(N)	
		$Z^4$ , CRTC(I)	X4(N)	
		$Z^5$ , RHO(I)	X5(N)	
		Zone masses, G(I)	EC(N)	
		Mass depletion energies, RFF(I)	PB(N)	
		Mass depletion area, SMLR(I)**2	SMLQ(N)	
		Left boundary of rezone region, R(IL)	RR	/NAMEC/ or internal
		Right boundary of rezone region, R(IR)	SAVRV6	
		Pusher velocity RD(1) or RD(IG-1)	RLF	
		Solid-vapor interface (see below)	RPUSH	
		LTE-non-LTE interface R(NR)	RNR	
		Left boundary of active region R(ICA)	RICA	
		Right boundary of active region R(ICB)	RICB	

The solid-vapor interface is defined by  $R(IA)$  for a left pusher (solid region) or  $R(IB)$  for a right pusher, provided that the vaporization flag  $BOILA$  is zero. If  $BOILA \neq 0$ ,  $RPUSH$  is located within the vaporizing zone at a position defined by the fraction of the mass in the zone which is still unvaporized. When the new zone structure is established, these mass fractions are recomputed if necessary.

The subroutine  $GENRAT$  is then called to fill in any previously unspecified quantities  $VAL(2)$ ,  $(3)$ ,  $(4)$ ,  $(5)$ , and subroutine  $MOVE$  is called to shift blank  $COMMON$  array elements with indices  $I \geq IR$  to their new locations. The rezone index  $IR$  for the current rezone region and the indices  $IL$  and  $IR$  in the  $RK$  array for yet unprocessed regions to the right of the current region are also shifted. Other indices and quantities which have been saved as described above are shifted to their new values. The  $ZONE$  routine (see below) generates new boundary positions, masses, and specific volumes in the rezone region. Boundary velocities are recomputed by linear spatial interpolation.

The next step is to recompute, for the new zones, the quantities  $E(I)$ ,  $FEW(I)$ ,  $RDD(I)$ ,  $SMLR(I)$ , and the five elemental mean ionic charges. This is done by mass-weighting (except for  $RDD$ ) to produce corresponding extensive quantities, computing the amount of each of these quantities in each subzone (subzones are defined by the union of the old and the new zone boundaries), assembling the subzone quantities belonging to each of the new zones, and dividing by the new zone masses (except for  $RDD$ ) to restore the original physical dimensions. Subroutine  $UP37A$  is called to update lambda-region arrays and clear the "save" locations of the old information which is no longer needed.

Next, subroutine "PROFL" (see below) is called if its flag was set, to eliminate by a special procedure any plateaus in the specific energy distribution in the solid which result from zone splitting. Finally, for LTE zones, the temperature and all of the thermodynamic variables except the specific internal energy are evaluated by an inversion of the state equation.

If the material identification number is in the range 201 to 300, this requires an iterative procedure, using interval halving or similar techniques. For non-LTE zones and LTE zones which specify the general ionic state equation routine EIONX (by material identification numbers  $\leq 200$  or  $> 300$ ), a more efficient procedure is being developed which makes use of the established elemental mean ionic charges and requires no iteration. At present, a simple version of this new procedure is used for non-LTE zones in one-material problems.

### 1.3. CONVER, ZONE, ZRAD, AND ZMS

The most important mathematical problem encountered in generating or rezoning a problem is the conversion between spatial coordinates and mass coordinates. The relationship between these is specified by a density profile, which is just a list of material densities for each of the original zones (in the rezoning case), or is given by input in generating a new problem. The four routines described in this section are the ones involving this conversion process, and are therefore discussed together.

Zoning may be carried out either with space or mass as the basic independent parameter. If VAL(6), the mass-space switch, is 1., the zones are to be set up with a uniform ratio VAL(4) of spatial widths, while VAL(6)=2. specifies a uniform ratio VAL(4) of zone masses. With space zoning, the conversion problem is one of integrating the density profile to obtain new zone masses, whereas with mass zoning, the problem is to integrate the reciprocal of the density profile to obtain new zone widths. The problem is nontrivial for three reasons:

1. The space variable may be the radius variable of a cylindrical or spherical coordinate system, in which case the interpretation to be given the density profile is not without ambiguity.
2. The zone widths must, in many cases, add up exactly to a value (the total region width) prescribed in advance, and the new boundary positions must conform to the zone widths as well as

possible within the limits of the finite computer word size. The sequences of boundary positions and zone widths need not be strictly derivable from each other, but they must both be regular and monotonic, with initial and terminal terms fixed. In some applications the boundary coordinates and zone widths are 8 to 10 orders of magnitude apart, and numerical methods designed for easier problems do not work.

3. For reasons of efficiency, neither involved iterative procedures nor indiscriminate use of double-precision operations is an acceptable approach to these problems.

The interpretation given the density profile in generating a new problem is that the complete density function is to be obtained by interpolating between the given points, and that the interpolant is linear not in the coordinate  $r$ , but in  $r^\alpha$  where  $\alpha=1,2,3$  for plane, cylindrical, and spherical symmetry respectively. This eliminates the need for solving anything more complicated than a quadratic equation in finding the width of a zone given the mass. In rezoning, the density profile is in origin a discrete mass profile, and any interpolation at all would be in a sense gratuitous. The assumption which is made throughout SPUTTER is that the density is constant throughout each zone; in rezoning, therefore, the profile is considered as defining a histogram, and the integration process is thus always linear. The integration subroutines ZMS and ZRAD are directed to these two alternative interpretations of the density profile by a switch IZSWT which is set in MAIN1 to -1 for rezoning and to 0 for generating.

The union of the set of new zone boundaries and the set of old zone boundaries or profile points thus defines a set of subzones within each of which the density is either constant or linear in  $r^\alpha$ . The mass integration routine ZMS is called for each subzone to determine the mass contained therein. The inverse problem is solved by the ZRAD subroutine, in which

each subzone has a given mass and it is required to find the volume from the profile. In these routines the evaluation of the slope from the profile is done in single precision, while the integration calculation itself may be done in double precision.

The CONVER routine is called from GENRAT to convert specifications into consistent form. Thus, if VAL(6)=2. specifying mass zoning, but any of VAL(1), VAL(2), or VAL(3) are specified with negative sign to indicate that they are widths, not masses, then the CONVER routine, with the aid of ZMS and the density profile, computes the proper masses where necessary for VAL(1) (region mass), VAL(2) (first zone mass), and VAL(3) (last zone mass). Similarly, CONVER calls ZRAD in the less usual case where space zoning is specified in terms of the mass of the region or of the first or last zones.

The ZONE routine computes the mass, width, boundary coordinates, and specific volume of each zone. After generating the zones using VAL(4) as the generating ratio, the subzone partition is established and ZMS (for space zoning) or ZRAD (for mass zoning) is called for each in turn to complete the description of the subzone. The assemblage of the subzones into the new zones to which they belong may be done with double numerical precision. The largest discrepancies arise for the boundary positions, which are cumulative. If the right region boundary is prescribed (always the case in rezoning and often the case in generating), the generated region will exceed or fall short of the target by an amount which is not necessarily small compared with the width of the last zone, even if double precision operations as described above are used. Smaller discrepancies arise for the width or mass of the last zone, since the prescribed value VAL(3) will not, in general, agree to the last digit with the value resulting from starting with VAL(2) and multiplying N times by VAL(4).

In each of these cases the discrepancies are removed by the following Procrustean device: The quantity in question is regenerated in reverse order (from right to left), using the prescribed right-end value as the initial condition. In calculating widths or masses, the initial condition is thus

VAL(3) and the generating ratio is  $1./VAL(4)$ , while in calculating boundary positions, the initial value is  $R(IR)$  and the increment is  $-DELTAR(I)$ , where  $DELTAR(I)$  is the same as the value used in the forward sweep. Each quantity obtained in the forward sweep is then multiplied by a factor  $W(I)$  and each reverse-sweep quantity by  $1.-W(I)$ , and the two results are then added to give the final result.  $W(I)$  is defined as that linear function of the index  $I$  which assumes the value 1. at the left end of the region and 0. at the right end.

It may be the case that this last procedure eliminates the need for double precision arithmetic at any stage of the calculation. Experimentation on this point is planned.

#### 1.4. PROFL

The PROFL subroutine is designed to eliminate a severe perturbation of the solid heat conduction calculation which results from splitting a single large zone into 10 or 20 small zones, each with the same specific internal energy and therefore the same temperature. The effect of this is to multiply the conductive flux at the ends of this region by a large factor and to destroy it entirely in the interior.

It is assumed that the conduction front is moving to the left, so that the temperature is a nondecreasing function of position  $r$ . The profile to be established is required to assume the same values at the ends of the region as the adjacent zones outside, and to conserve the original total energy content. These three conditions are sufficient to determine a quadratic function but not to ensure that the quadratic is a non-decreasing function of  $r$ . The profile is therefore defined as a quadratic which, starting at the right boundary, decreases monotonically, reaching the left boundary value at a point which may be in the interior. The profile in that case is continued as a constant to the left boundary. The energy in each zone is then simply the integral of this profile across the zone.

The geometry may have plane, cylindrical, or spherical symmetry.

## 1.5. APPENDIX: CODE LISTINGS

```

SUBROUTINE AUT0RZ
C
C   COMPILLED MAY 16, 1966
C
C*****C*****
C
C*          S P U T T L E   C O M P O U
C*
C*
COMMON  LMDA(37), NR      , USMR , IA      , IB      , ICA      , ICH      ,
1  KMAX  , BLANK1, BLANK2, BLANK3, IAP1  , IBP1  , ICAP1  , ICBP1  ,
2  II    , IG      , NRAD  , BLANK4, IAM1  , IBM1  , ICAM1  , ICBM1  ,
3  IIP1  , IGM1  , IALPHA, BLANK5, TH      , TMAX  , BLANK6, DELPRT,
4  FREQ  , CNTMAX, AR      , ASMR , PUSHA , PUSHB , BOTIA , BOTLB ,
5  CVA   , CVB   , SLUG  , ALPHA , HVA   , HVB   , HCA   , HCB   ,
6  EMINA , EMINB , CA     , CB     , GA     , GB     , GL     , GR     ,
7  RHOL  , RHOR  , EP10  , LPS1   , RIA    , RIB    , RDTA   , RDTB   ,
8  RPIA  , RPIB  , RPDIA , RPDIB , TRINT , IA     , TH     , IC     ,
COMMON  TD      , TE      , DTH2  , DTH2P , DTH1  , DTHM1 , DTHAX ,
1  DTMAX1, DTMAX2, DTMAX3, DTR    , SWITCH, CO     , CMTH  , DELTA ,
2  GAMA  , WCRIT , SIGMAQ, AC      , ACOSTH, CHVRT , SUMPA , SUMRB ,
3  ROIA  , ROIAM1, ROIB   , ROIBP1, GMS   , S1     , S2     , S3     ,
4  S4     , S5     , S6     , S7     , S8     , S9     , S10    , S11    ,
5  S12    , S13    , S14    , S15    , S16    , S17    , S18    , S19    ,
6  S20    , EO     , FO     , TAU   , ZERO  , R      (152), DELTAR(152),
7  ASQ    (152), RD      (152), VD      (152), RDD    (152), SMLR   (152),
8  DELR   ( 37), P      (152), P1     (152), PH      (152), PB1    (152),
COMMON  P2     (152), SV      (152), ITHO   (152), THETA (152),
1  W      (152), E      (152), EI      (152), EK      (152), A      (152),
2  V      (152), G      (152), D      (152), C      (152), X2     (152),
3  X3     (152), X4     (152), X5     (152), X6     (152), X7     (152),
4  SMLA   (152), SMLB   (152), SMLC   (152), SMLD   (152), SMLE   (152),
5  EC     (152), EK     (152), SMLG   (152), SMLH   (152), BIGA   (152),
6  BIGB   (152), CV     (152), BC     (152), BR     (152), CHIC   (152),
7  CHIR   (152), CAPAC  (152), CAPAR  (152), CRTC   (152), CRTK   (152),
8  CRTPC  (152), GOFB   (152), FEW    (152), CAR    (152), OKLM   ( 37),
COMMON  TELM   ( 37), EKLK   ( 37), ELM    ( 37), FCLM   ( 37),
1  FRLM   ( 37), WLM    ( 37), OLM    ( 37), AMASNO ( 37), CHRNO  ( 37),
2  ZP1    ( 37), ZP2    ( 37), SOLID  ( 37), ECHCK  ( 37), RK     (104),
3  RL     ( 37), RHOK   (104), RDK    (104), THETAK(104), TEMP   ( 16),
4  HEAD   ( 12), MAXL    , MAXLM
C
C*****C*****
C
C*          INPUT DATA *****
C
C   THETAK(53)=SOLID REZONE RATIO
C   THETAK(54)=SOLID REZONE INTERFACE
C   THETAK(55)=NUMBER OF SMALL VAPOR ZONES
C   THETAK(56)=SSIGMA IN VAPOR TO RIGHT OF INTERFACE
C   THETAK(57)=REZONE CRITERION
C   THETAK(58) = 1. FOR SOLID REZONE ONLY, = 2. FOR SOLID AND VAPOR
C   THETAK(58) .EQ.-1. FOR REZONE IN VAPOR ONLY
C   THETAK(59) = RT. BOUNDARY FOR CONSTANT ZONES
C   THETAK(60) = CONSTANT ZONE SIZE
C   IF (THETAK(58).EQ.0.) GO TO 1000
C   ISR=THETAK(54)
C   NREZ=THETAK(57)
C   NSVZ = THETAK(55)
C   IF (IA.LE.ISR+NREZ) GO TO 10
C   IF (THETAK(58).EQ.1.) GO TO 1000
C   IF (THETAK(59).EQ.0.) GO TO 1000
C   IF (R(16).LT.THETAK(59)) GO TO 1000
C   RZCF = 1.
C   GO TO 11
10  RZCF = 0.
C   THETAK(54)=ISR-1

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      IF (THETAK(58).EQ.1.) GO TO 100
      IF (THETAK(59).EQ.0.) GO TO 10
11  K = 0
      DO 12 I = IA, IB
      J = IA + IB - I
      IF (R(J).GT.THETAK(59)) GO TO 12
      IF (K.EQ.0) K = J
      IF (DELTA(J).LT.THETAK(60)) GO TO 13
12  CONTINUE
C    K = RT. INDEX FOR CONSTANT ZONES
C    J = LF. INDEX FOR CONSTANT ZONES
C    - - FULL REZONE - -
13  IF (RZCF.GT.0. .AND. DELTA(K).LT.1.5*THETAK(60)) GO TO 1000
      IF (RZCF.EQ.0. .AND. K-J.LT.3) GO TO 20
      RK(2) = J + 1
      RK(3) = K + 1
      RK(4) = -THETAK(60)
      RK(5) = 0.
      RK(6) = 0.
      RK(7) = INT((R(K+1)-R(J)) / THETAK(60)) + 1
      RK(8) = 1.
      IF (IB.GE.(K + 1 + NSVZ)) GO TO 14
      IF (RZCF.EQ.1.) GO TO 499
      GO TO 21
14  RK(13) = K + 1
      RK(14) = IB
      RK(15) = G(K)
      RK(16) = 0.
      RK(17) = THETAK(56)
      RK(18) = 0.
      RK(19) = 2.0
      IF (RZCF.EQ.1.) GO TO 40
      RK(24) = THETAK(54)
      RK(25) = -ISR
      RK(26) = G(ISR-1) / THETAK(53)
      RK(27) = RK(26)
      RK(28) = 1.0
      RK(29) = THETAK(53)
      RK(30) = 2.0
      RK(1) = 3.
      GO TO 500
C    - - OUTER VAPOR FIRST - -
20  IL = IA + NSVZ
      IF (THETAK(59).NE.0.) IL = K
      IF (IB.LE.(IL+NSVZ)) GO TO 100
      RK(2) = IL
      RK(3) = IB
      RK(4) = G(IL-1)
      RK(5) = 0.
      RK(6) = THETAK(56)
      RK(7) = 0.
      RK(8) = 2.0
      AUT00700
      AUT00720
      AUT00730
      AUT00740
      AUT00750
      AUT00760
21  IF (THETAK(58).EQ.(-1.)) GO TO 499
C    - - SOLID AFTER VALOR - -
30  RK(13) = THETAK(54)
      RK(14) = -ISR
      RK(15) = G(ISR-1)/THETAK(53)
      RK(16) = RK(15)
      RK(17) = 1.0
      RK(18) = THETAK(53)
      RK(19) = 2.0
      AUT00620
      AUT00650
      AUT00670
40  RK(1) = 2.
      GO TO 500
      AUT00770
C    - - SOLID ONLY - -
100 IF (THETAK(58).EQ.(-1.)) GO TO 1000
      IF (IB+INT(THETAK(53)).GT.150) GO TO 200
      RK(2) = THETAK(54)
      RK(3) = -ISR
      AUT00790

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      RK(4) = G(ISR-1)/THETAK(SJ)
      RK(5) = RK(4)
      RK(6) = 1.
      RK(7) = THETAK(SJ)
      RK(8) = 2.0
499  RK(1) = 1.
500  S16 = -1.
      CALL MPI
200  S1 = 2.0200
      CALL UNCLE
1000 RETURN
      END

```

AUT00810

AUT00840

AUT00850

AUT00870

AUT00880

AUT00890

AUT00900

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SUBROUTINE REZONE
COMPILED MAY 16, 1966

SUBROUTINE REZONE CALCULATES NEW ZONAL PROPERTIES
BETWEEN ANY TWO POINTS (IL,IR), SHIFTS THE PRESENT
ARRAYS TO FIT THESE POINTS.
CALLED FROM EITHER IMAIN OR AUTOZ

      S P U T T E R   C O M M O N

COMMON  LMDA(37), NR      , NSMLR , IA      , IB      , ICA      , ICB      ,
1  KMAX , BLANK1, BLANK2, BLANK3, IAP1 , IBP1 , ICAP1 , ICBP1 ,
2  II   , IG      , NRAD , BLANK4, IAM1 , IBM1 , ICAM1 , ICBM1 ,
3  IIP1 , IGM1 , IALPHA, BLANK5, TH      , TMAX , BLANK6, DELPRT,
4  FREQ , CNTMAX, AR      , ASMLR , PUSHA , PUSHB , BOILA , BOILB ,
5  CVA   , CVB   , SLUG , ALPHA , HVA   , HVB   , HCA   , HCB   ,
6  EMINA , EMINB , CA      , CB      , GA      , GB      , GL      , GR      ,
7  RHOL , RHOR , EPIO , EPSI , RIA      , RIB      , RDIA , RDIB ,
8  RPIA , RPIB , RPDIA , RPDIB , TPRINT, TA      , TB      , TC      ,
COMMON  TD      , TE      , DTH2 , DTH2P , DTH1 , DTRMIN, DTMAX ,
1  DTMAX1, DTMAX2, DTMAX3, DTR , SWITCH, CO      , CMIN , DELTA ,
2  GAMA , WCRIT , SIGMAQ, AC      , ACO3T4, CNVRT , SUMRA , SUMRB ,
3  ROIA , ROIAM1, ROIB , ROIBP1, GMS , S1      , S2      , S3      ,
4  S4      , S5      , S6      , S7      , S8      , S9      , S10     , S11     ,
5  S12     , S13     , S14     , S15     , S16     , S17     , S18     , S19     ,
6  S20     , E0      , F0      , TAU , ZERO , R      (152), DELTAR(152),
7  ASQ      (152), RD      (152), VD      (152), RDD      (152), SMLR      (152),
8  DELR      ( 37), P      (152), P1      (152), PB      (152), PB1      (152),
COMMON  P2      (152), SV      (152), RHO      (152), THETA (152),
1  W      (152), E      (152), EI      (152), EK      (152), A      (152),
2  V      (152), G      (152), D      (152), C      (152), X2      (152),
3  X3      (152), X4      (152), X5      (152), X6      (152), X7      (152),
4  SMLA      (152), SMLB      (152), SMLC      (152), SMLD      (152), SMLE      (152),
5  EC      (152), ER      (152), SMLQ      (152), SMLH      (152), BIGA      (152),
6  BIGB      (152), CV      (152), BC      (152), BR      (152), CHIC      (152),
7  CHIR      (152), CAPAC      (152), CAPAR      (152), CRTC      (152), CRTR      (152),
8  CRTPC      (152), GOFR      (152), FEW      (152), CAR      (152), OKLM      ( 37),
COMMON  TELM      ( 37), EKLM      ( 37), ELM      ( 37), FCLM      ( 37),
1  FRLM      ( 37), WLM      ( 37), OLM      ( 37), AMASNO( 37), CHRNO      ( 37),
2  ZP1      ( 37), ZP2      ( 37), SOLID      ( 37), ECHCK      ( 37), RK      (104),
3  RL      ( 37), RHOK      (104), RDK      (104), THETAK(104), TEMP      ( 16),
4  HEAD      ( 12), MAXL      , MAXLM

*****
COMMON /NAMEC/      RLF , RINC , RPUSH , RR      , DR      ,
COMMON /NAMEC/      KCNT , LCNT , MCNT , N      , INO , IL , IR
COMMON /NAMEC/ VAL      ( 10), RTABRO(150), RTABTH(150), RTABRD(150),
1  RHOTAB(150), THETAB(150), RDTAB(150)
COMMON /PROSW/ IPROSW
COMMON /LMSB/ U(1) /LMSC/ M(1)

DIMENSION XZONE(11,9)
DIMENSION ZN(6)
DIMENSION ZK(5), Z(1), WT(1)
EQUIVALENCE (M(1),KMX), (M(2),Z(1)), (M(3),WT(1))
EQUIVALENCE (RK(2),XZONE)
REAL 1K
LOGICAL IBOIL

SET THE NUMBER OF REGIONS TO BE CHANGED

NCARD = RK(1)

```

C	DO 140 J = 1,NCARD	REZ00680
C		REZ00690
C	INITIALIZE VARIABLES THAT KEEP TAB ON AREAS TO BE REZONED	REZ00700
C		REZ00710
	IRN = 0	REZ00720
	IRO = 0	REZ00730
	IGO = IG	REZ00740
	IL = XZONE(1,J)	REZ00750
	IR = XZONE(2,J)	REZ00760
C		REZ00770
C	DO WE TURN ON THE PROFILE SWITCH	REZ00780
C		REZ00790
	IPROSW = 0	REZ00800
	IF (IR .LT. 0) IPROSW = 1	REZ00810
	IR = IABS(IR)	REZ00820
	SAVRV6 = R(IR)	REZ00840
C		REZ00850
C	SAVE THE APPROPRIATE ARRAYS	REZ00860
C		REZ00870
	N = 1	REZ00890
C		REZ00900
	DO 99 I = IL,IR	REZ00910
C		REZ00920
C	SAVE TABLES	REZ00930
C		REZ00940
	RDTAB(N) = RD(I)	
	RTABRD(N) = R(I)	
	IF (I.EQ.IR) GO TO 99	
C		REZ00970
	RHOTAB(N) = 1./SV(I)	REZ00980
C		REZ00990
	RTABRO(N) = R(I)	REZ01000
	THETAB(N) = E(I)	REZ01010
	RTABTH(N) = R(I)	REZ01020
	CAR(N)=FEW(I)	REZ01030
	CAPAC(N)=X6(I)	REZ01040
	CHIC(N)=BC(I)	REZ01050
	CHIR(N)=BR(I)	REZ01060
	X4(N)=CRTC(I)	REZ01070
	X5(N)=RHO(I)	REZ01080
	EC(N)=G(I)	REZ01090
	IF (EC(N).LE.0.) GO TO 110	
	PB(N)=RDD(I)	REZ01100
	SMLQ(N)=SMLR(I)**2	REZ01110
C		REZ01120
	N = N + 1	REZ01140
	IF (N + 1.GT.150) GO TO 110	REZ01150
	99 CONTINUE	REZ01160
C		REZ01170
	KCNT = IR - IL	
	LCNT = KCNT	
	MCNT = KCNT + 1	REZ01190
C		REZ01210
	IF(PUSHA) 104,106,105	REZ01220
104	RLF = RD(IG-1)	REZ01230
	GO TO 111	REZ01240
105	RLF = RD(1)	REZ01250
106	GO TO 111	REZ01260
C		REZ01270
110	S1 = 55.0110	REZ01280
	CALL UNCLE	REZ01290
C		REZ01300
111	IBOIL = .FALSE.	REZ01310
	IF(PUSHA) 112,115,113	REZ01320
112	RPUSH = R(IB)	REZ01330

IF(IL.GT.IB .OR. IR.LT.IB .OR. HOILA.EQ.0.) GO TO 115	REZ01340
IBB = IB	REZ01350
JB1 = IBP1	REZ01360
X = -1.	REZ01370
GO TO 114	REZ01380
113 RPUSH = R(IA)	REZ01390
IF(IL.GT.IAM1.OR.IR.LT.IAM1 .OR. HOILA.EQ.0.) GO TO 115	REZ01400
IBB = IAM1	REZ01410
JB1 = IAM1	REZ01420
X = 1.	REZ01430
114 IBOIL = .TRUE.	REZ01440
RPUSH=R(JB1)+X*(1.-SOLID(24))*G(IBM)*SOLID(17)/A(JB1)	REZ01450
115 CONTINUE	REZ01460
C	REZ01470
C	REZ01480
C	REZ01490
IF((IR-IL) .GT. 149) GO TO 100	REZ01500
C	REZ01510
C	REZ01520
C	REZ01530
VAL(1) = 0.	REZ01540
DO 120 I = 2,10	REZ01550
VAL(I) = XZONE(I+1,J)	REZ01560
120 CONTINUE	REZ01570
IRM1 = IR - 1	REZ01580
IF(VAL(6) .EQ. 2.) GO TO 801	REZ01590
IF(VAL(6) .NE. 1.) GO TO 100	REZ01600
DO 800 I = IL,IRM1	REZ01610
800 VAL(1) = VAL(1) - DELTAR(I)	REZ01620
GO TO 803	REZ01630
801 DO 802 I = IL,IRM1	REZ01640
802 VAL(1) = VAL(1) + G(I)	REZ01650
803 CONTINUE	REZ01660
RR = R(IL)	REZ01670
CALL GENRAT	REZ01680
C	REZ01690
129 CONTINUE	REZ01700
HNR = R(NR)	REZ01710
RICA = R(ICA)	REZ01720
RICB = R(ICB)	REZ01730
C	REZ01740
C	REZ01750
C	REZ01760
SHIFT ARRAYS FOR NEW CONDITIONS	REZ01770
CALL MOVE	REZ01780
C	REZ01790
C	REZ01800
C	REZ01810
ADJUST BOUNDARIES	REZ01820
IRO = IR	REZ01830
IRN = IL + IFIX(VAL(5))	REZ01840
IR = IRN	
ITEM = IRN-IRO	
IRM1 = IR - 1	
R(IR) = SAVRV6	
C	REZ01850
C	REZ01860
C	REZ01870
CALL ZONE	REZ01880
C	REZ01890
C	REZ01900
C	REZ01910
DO 1294 I = IL,IRM1	REZ01920
IF(S10) 1291,1292,1293	REZ01930
1291 A(I) = 1.	REZ01940
GO TO 1294	REZ01950
1292 A(I) = 2. * R(I)	REZ01960

GO TO 1294	
1293 A(I) = 3. * R(I)**2	REZ01970
1294 CONTINUE	REZ01980
C	REZ01990
IF (J.EQ.NCARD) GO TO 1300	REZ02000
JP1 = J + 1	
DO 130 I = JP1,NCARD	
IF (XZONE(1,J).GT.XZONE(2,I)) GO TO 130	
XZONE(1,I) = XZONE(1,I) + FLOAT(ITEM)	REZ02030
XZONE(2,I) = XZONE(2,I) + FLOAT(ITEM) * SIGN(1.,XZONE(2,I))	REZ02040
130 CONTINUE	REZ02050
1300 CONTINUE	
C	REZ02060
IF(PUSHA) 136, 137, 131	REZ02070
131 CONTINUE	REZ02080
IB = IB+ITEM	REZ02090
IG = IG+ITEM	REZ02100
ICB = IB	REZ02110
IF(IA.GT.IL) IA = ISAV1(IG,IA,RPUSH,R(1))	REZ02120
IF(ICA.GT.IL) ICA = ISAV1(IG,-ICA,RICA,R(1))	REZ02130
NR = ISAV1(IG,0,RNR,R(1))	
GO TO 139	REZ02150
C	REZ02160
136 CONTINUE	REZ02170
IG=IG+ITEM	REZ02180
IF(IB.GT.IL) IB = ISAV1(IG,-IB,RPUSH,R(1))	REZ02190
ICB=ISAV1(IG,ICB,RICB,R(1))	REZ02200
NR = ISAV1(IG,0,RNR,R(1))	
GO TO 139	REZ02220
C	REZ02230
137 CONTINUE	REZ02240
IG = IG+ITEM	REZ02250
IB = IB + ITEM	REZ02260
ICB = ISAV1(IG,ICB,RICB,R(1))	REZ02270
IF(ICA.GT.IL) ICA = ISAV1(IG,-ICA,RICA,R(1))	REZ02280
NR = ISAV1(IG,0,RNR,R(1))	
139 CONTINUE	REZ02300
C	REZ02310
IAP1 = IA+1	REZ02320
IAM1 = IA-1	REZ02330
IBP1 = IB+1	REZ02340
IBM1 = IB-1	REZ02350
ICAM1 = ICA-1	REZ02360
ICAP1 = ICA+1	REZ02370
ICBM1 = ICB - 1	REZ02380
ICBP1 = ICB + 1	REZ02390
IGM1 = IG-1	REZ02400
C	REZ02410
C	REZ02420
C	REZ02430
UPDATE VAPOR FRACTION OF BOILING ZONE	REZ02440
IF(.NOT.IBOIL) GO TO 1398	REZ02450
IF(PUSHA) 1392,1398,1394	REZ02460
1392 IBB = IB	REZ02470
JB1 = IBP1	REZ02480
GO TO 1396	REZ02490
1394 IBB = IAM1	REZ02500
JB1 = IAM1	REZ02510
1396 SOLID(24)=1.-(RPUSH-R(JB1))*A(JB1)*X/(G(IBB)*SOLID(17))	REZ02520
1398 CONTINUE	REZ02530
C	REZ02540
C	REZ02550
C	REZ02560
CALCULATE NEW VELOCITY	REZ02570
DO 200 I=IL,IR	REZ02580
RDOT=R(I)	
LASTN=3	

CALL INTERP(RDOT, LASTN)	REZ02590
RU(1)=RDOT	REZ02600
200 CONTINUE	REZ02610
C	REZ02620
C	REZ02630
C	REZ02640
C	REZ02650
BALANCE THE NEW ENERGIES	REZ02660
REGENERATE ZONE ENERGY, MASS DEPLETION AND NONEQ VARIABLES	REZ02670
K=1	REZ02680
SM1=EC(1)	REZ02690
I=IL	REZ02700
SM2=G(1L)	REZ02710
300 EN=0.	REZ02720
IF(SUMRA.EQ.0.) GO TO 305	REZ02730
RN=0.	REZ02740
SN=0.	REZ02750
305 IF(ZP1(26).EQ.0.) GO TO 315	REZ02760
DO 310 JJ = 1,6	REZ02770
ZN(JJ) = 0.	REZ02780
310 CONTINUE	REZ02790
315 SM=AMIN1(SM1,SM2)	REZ02800
EN=EN+THETAB(K)*SM	REZ02810
IF(SUMRA.EQ.0.) GO TO 320	REZ02820
RN=RN+PB(K)*SM/EC(K)	REZ02830
SN=SN+SMLQ(K)*SM	REZ02840
320 IF(ZP1(26).EQ.0.) GO TO 325	
ZN(1)=ZN(1)+CAR(K)*SM	
IF (PUSHA.GE.0..AND.RTABRO(K).LT.RNR) GO TO 325	
IF (PUSHA.LT.0..AND.RTABRO(K).GE.RNR) GO TO 325	
ZN(2)=ZN(2)+CAPAC(K)*SM	REZ02850
ZN(3)=ZN(3)+CHIC(K)*SM	REZ02860
ZN(4)=ZN(4)+CHIR(K)*SM	REZ02870
ZN(5)=ZN(5)+X4(K)*SM	REZ02880
ZN(6)=ZN(6)+X5(K)*SM	REZ02890
325 SM1=SM1-SM	REZ02900
SM2=SM2-SM	REZ02910
IF(SM1.GT.0.) GO TO 330	REZ02920
IF (K.LT.KCNT) K = K + 1	
SM1=EC(K)	
330 IF(SM2.GT.0.) GO TO 315	REZ02940
E(I)=EN/G(I)	REZ02950
E1(I)=E(I)	REZ02960
IF(SUMRA.EQ.0.) GO TO 335	REZ02970
RDD(I)=RN	REZ02980
SMLR(I)=SQRT(ABS(SN/G(I)))	REZ02990
335 IF(ZP1(26).EQ.0.) GO TO 340	REZ03000
IF (PUSHA.GE.0..AND.I.LT.NR) GO TO 340	REZ03010
IF (PUSHA.LT.0..AND.I.GT.NR-1) GO TO 340	
FEW(I)=ZN(1)/G(I)	REZ03020
X6(I)=ZN(2)/G(I)	REZ03030
BC(I)=ZN(3)/G(I)	REZ03040
BR(I)=ZN(4)/G(I)	REZ03050
CRTC(I)=ZN(5)/G(I)	REZ03060
RHO(I)=ZN(6)/G(I)	REZ03070
SUMZ=X6(I)*WT(1)	
IF(KMX.EQ.1) GO TO 337	
SUMZ=SUMZ + BC(I)*WT(11)	
IF(KMX.EQ.2) GO TO 337	
SUMZ=SUMZ + BR(I)*WT(21)	
IF(KMX.EQ.3) GO TO 337	
SUMZ=SUMZ + CRTC(I)*WT(31)	
IF(KMX.EQ.4) GO TO 337	
SUMZ=SUMZ + RHO(I)*WT(41)	
337 IF (SUMZ.EQ.0.) GO TO 340	
X6(I) = X6(I) * FEW(I) / SUMZ	
IF (KMX.EQ.1) GO TO 338	



```

      BC(I) = BC(I) * FEW(I) / SUMZ
      IF (KMX.EQ.2) GO TO 338
      BR(I) = BR(I) * FEW(I) / SUMZ
      IF (KMX.EQ.3) GO TO 338
      CRTC(I) = CRTC(I) * FEW(I) / SUMZ
      IF (KMX.EQ.4) GO TO 338
      RHO(I) = RHO(I) * FEW(I) / SUMZ
338  CONTINUE
340  I=I+1
      SM2=G(I)
      IF(I.LT.IR) GO TO 300
      CALL UP37A(IRO,IGO)
C
C    CALCULATE ZONE TEMPERATURES
C
      IF(IPROSW.EQ.1) CALL PROFL
      DO 380 I=IL,IRM1
      IF (ZP1(26).EQ.0.) GO TO 342
      IF (PUSHA.GE.0..AND.I.GE.NR) GO TO 364
      IF (PUSHA.LT.0..AND.I.LT.NR) GO TO 364
342  IC = 0
      TL=0.
      TR=2.*THETA(I-1)
      IF(I.EQ.IL)TR=2.*THETA(I)
345  THETA(I)=.5*(TL+TR)
      SMLA(I) = THETA(I)
      CALL EOS(I)
      IF(IC.GT.30)GO TO 365
      DE=E(I)-EI(I)
      IF(ABS(DE/EI(I)).LE.S19) GO TO 370
      IF(DE.GT.0.) GO TO 350
      IF(IC.GT.0) GO TO 355
      TL=THETA(I)
      TR=TL*2.
      GO TO 345
350  TR=THETA(I)
      GO TO 360
355  TL=THETA(I)
360  IC=IC+1
      GO TO 345
365  WRITE(6,1000) I
1000  FORMAT(29H0ENERGY DISCONTINUOUS IN ZONE,13)
      GO TO 380
364  ZK(1) = X6(I)
      ZK(2) = BC(I)
      ZK(3) = BR(I)
      ZK(4) = CRTC(I)
      ZK(5) = RHO(I)
      EION = 0.
      DO 368 K = 1,KMX
      KK = 10*K - 9
      LMAX = Z(KK)
      IK = AINT(ZK(K))
      L = INT(IK)
      NV = M(KK+3) + 1
      IF (L.EQ.0) GO TO 367
      DO 366 JJ = 1,L
      NV = NV + 1
      EION = EION + WT(KK) * U(NV)
366  CONTINUE
      IF (L.EQ.LMAX) GO TO 368
367  EION = EION + WT(KK)*U(NV+1)*(ZK(K)-IK)
368  CONTINUE
369  AMASNI = AMASN(I)
      CV(I) = 1.5 * AMASNI * (1.0 + FEW(I))

```

REZ03120  
 REZ03130  
 REZ03140  
 REZ03460  
 REZ03150  
 REZ03160  
 REZ03170  
 REZ03180  
 REZ03190

REZ03210  
 REZ03220  
 REZ03230  
 REZ03240

REZ03250  
 REZ03260  
 REZ03270  
 REZ03280  
 REZ03290  
 REZ03300  
 REZ03310  
 REZ03320  
 REZ03330  
 REZ03340  
 REZ03350  
 REZ03360  
 REZ03370  
 REZ03380  
 REZ03390  
 REZ03400  
 REZ03410

```

      THETA(I) = (E(I) - AMASNI*EION - EZERO(I)) / CV(I)
      P1(I) = AMASNI * (1.0 + FEW(I)) * THETA(I) / SV(I)
370  EI(I)=E(I)
380  CONTINUE
C
C
C
140  CONTINUE
      RETURN
C
C
      ERROR
100  S1 = 55.0100
      CALL UNCLE
C
      END

```

02/23/66

REZ03420  
 REZ03430  
 REZ03440  
 REZ03450  
 REZ03470  
 REZ03480  
 REZ03490  
 REZ03500  
 REZ03510  
 REZ03520  
 REZ03530  
 REZ03540  
 REZ03550



```

SUBROUTINE CONVR
COMPILED JUNE 6, 1966
C*****
C
C*          S P U T T E R   C O M M O N
C*
C*          COMMON LMDA(9405)
C*          DIMENSION EC(152)
C*          EQUIVALENCE (S1, LMDA(136)), (EC(1), LMDA(5517))
C
C*****
C*          COMMON /NAMEC/ RLF, RINC, RPUSH, RR, DR
C*          COMMON /NAMEC/ KCNT, LCNT, MCNT, N, INO, IL, IR
C*          COMMON /NAMEC/ VAL (10), RTABRO(150), RTABTH(150), RTABPD(150)
C*          1 RHOTAB(150), THETAB(150), RDTAB (150)
C
C*          COMMON /SHEZS/ IZSWT
C*          DOUBLE PRECISION R0,R1,R2,R3,G1,G2,V5,ZMS
C
C*          M   M1   M2   (SPACE COORDINATES GIVEN)
C*          X   X   X   SPECIAL CASE   (X MEANS VALUE GIVEN)
C*          X   X   0
C*          X   0   X   SPECIAL CASE
C*          0   X   X   ERROR
C*          0   0   X   ERROR
C*          0   X   0
C*          X   0   0
C
C*          MASS - SPACE SWITCH
C*          IV6 = VAL(6) + 0.1
C
C*          ITEM = 0
C*          TEM = ABS(VAL(1))-ABS(VAL(3))
C*          DO 190 I = 1,3
C*          IF((VAL(6)-1.5)*VAL(I).GE.0.) GO TO 190
C*          R0=DBLE(RR)
C*          K=MINO(KCNT,2)
C*          80 IF(RTABRO(K).GT.RR.OR.K.EQ.KCNT) GO TO 90
C*          K=K+1
C*          GO TO 80
C*          90 CONTINUE
C*          GO TO (110,150),IV6
C
C*          SPACE COORDINATES
C*          110 CONTINUE
C
C*          CHECK FOR MIS-SPECIFICATION
C*          IF(VAL(3).GT.0.0 .AND. VAL(1).EQ.0.0) GO TO 180
C
C*          CHECK FOR SPECIAL CASE
C*          IF(I.EQ.3 .AND. (VAL(1).GT.0.0 .AND. VAL(3).GT.0.0)) GO TO 130
C*          120 CONTINUE
C*          R1=R0
C*          IF(KCNT.GT.2) GO TO 122
C*          G1 = 1.D20
C*          GO TO 124
C*          122 IF(SNGL(R1).EQ.RTABRO(K-1).AND.IZSWT.EQ.(-1)) GO TO 123
C*          R2=DBLE(RTABRO(K))
C*          G1=ZMS(R1,R2,K)
C*          GO TO 124
C*          123 G1=DBLE(EC(K-1))
C*          124 G2=DBLE(VAL(I))
C*          IF(G2.GE.G1) GO TO 126

```

```

CONV0010
CONV0020
CONV0030
CONV0040
CONV0050
CONV0060
CONV0080
CONV0090
CONV0100
CONV0110
CONV0120
CONV0130
CONV0140
CONV0150
CONV0180
CONV0190
CONV0200
CONV0210
CONV0220
CONV0230
CONV0240
CONV0250
CONV0260
CONV0270
CONV0280
CONV0290
CONV0300
CONV0310
CONV0320
CONV0330
CONV0350
CONV0360
CONV0370
CONV0410
CONV0420
CONV0430
CONV0440
CONV0450
CONV0460
CONV0470
CONV0480
CONV0490
CONV0500

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AFWL-TR-66-108, Vol III

125	CALL ZRAD(R1,R3,G2,VS,K)	
	GO TO 128	
126	IF(G2.GT.G1) GO TO 127	
	R3=R2	
	GO TO 128	
127	R1=R2	
	G2=G2-G1	
	IF(K.EQ.KCNT) GO TO 125	
	K=K+1	
	GO TO 122	
128	VAL(I)=SNGL(R3-R0)	
	GO TO 190	
C		CONV0590
130	CONTINUE	CONV0600
	VAL(3) = TEM	CONV0610
	ITEM = 1	CONV0620
	GO TO 120	CONV0630
C		CONV0640
C	MASS COORDINATES	CONV0650
C		CONV0660
150	CONTINUE	CONV0710
C		CONV0720
C	CHECK FOR MIS-SPECIFICATION	CONV0730
C		CONV0740
	IF(VAL(3).LT.0.0 .AND. VAL(1).EQ.0.0) GO TO 180	CONV0750
C		CONV0760
C	CHECK FOR SPECIAL CASE	CONV0770
C		CONV0780
	IF(1.EQ.3 .AND. (VAL(1).LT.0.0 .AND. VAL(3).LT.0.0)) GO TO 170	CONV0790
C		CONV0800
160	G1 = 0.00	
	RJ=R0+DBLE(ABS(VAL(I)))	
	R1=R0	
	IF(KCNT.GE.2) GO TO 162	
161	G1=ZMS(R1,R3,K)+G1	
	GO TO 165	
162	R2=DBLE(RTABRO(K))	
	IF(R2.GT.R3) GO TO 161	
	IF(SNGL(R1).EQ.RTABRO(K-1).AND.IZSWT.EQ.(-1)) GO TO 163	
	G1=ZMS(R1,R2,K)+G1	
	GO TO 164	
163	G1=DBLE(EC(K-1))+G1	
164	IF(R2.EQ.R3) GO TO 165	
	R1=R2	
	IF(K.EQ.KCNT) GO TO 161	
	K=K+1	
	GO TO 162	
165	VAL(I) = -SNGL(G1)	
	GO TO 190	
C		CONV0870
170	CONTINUE	CONV0880
	VAL(I) = TEM	CONV0890
	ITEM = 1	CONV0900
	GO TO 160	CONV0910
C		CONV0920
180	CONTINUE	CONV0930
181	S1 = 43.0181	CONV0940
	WRITE (6,10)	CONV0950
	RETURN	CONV0960
C		CONV0970
190	CONTINUE	CONV0980
	IF(ITEM.EQ.0) RETURN	CONV0990
	VAL(3) = ABS(VAL(1)) - ABS(VAL(3))	CONV1000
	RETURN	CONV1010
C		CONV1020
10	FORMAT ( 52H1A MIS-SPECIFICATION HAS BEEN FOUND IN THE ZONE DATA)	CONV1030
	END	CONV1040

```

SUBROUTINE ZONE
  COMPILED JULY 15, 1966
  *****
  S P U T T I R   C O M M O N
  *****

  COMMON  LMDA(37), NR      , NSMR , IA      , IB      , IC      , ICH
1  KMAX   , BLANK1, BLANK2, BLANK3, IAP1   , IBP1   , ICAP1 , ICHP1
2  II     , IG     , NRAD  , BLANK4, IAM1   , IBM1   , ICAM1 , ICBM1
3  IIP1   , IGM1   , IALPHA, BLANK5, TIM1   , TMAX   , BLANK6, DELPRT
4  FREQ   , CNTMAX, AK     , ASMR , PUSH1  , PUSH2  , BOT1A , BOT1B
5  CVA    , CVB    , SLUG  , ALPHA , HVA    , HVB    , HCA    , HCB
6  EMINA   , EMINB , CA     , CB     , GA     , GB     , GL     , GR
7  RHOL   , RHOR   , EPSI  , EPSI  , RIA    , RIB    , RIDIA , RIDIB
8  RPIA    , RPIB   , RPDIA , RPDIB , TPRINT, TA     , TB     , TC
  COMMON  TD      , TE      , DTHP , DTH2P , DTH3 , DTRMIN, DTMAX
1  DTMAX1, DTMAX2, DTMAX3, DTR   , SWITCH, CO     , CMIN  , DELTA
2  GAMA   , WCRIT , SIGMAQ, AC     , ACO3T4, CNVRT , SUMRA  , SUMRB
3  ROIA   , ROIAM1, ROIB   , ROIBP1, GMS    , S1     , S2     , S3
4  S4     , S5     , S6     , S7     , S8     , S9     , S10    , S11
5  S12    , S13    , S14    , S15    , S16    , S17    , S18    , S19
6  S20    , EO     , FO     , TAU    , ZERO   , R      (152), DELTAR(152)
7  ASQ    (152), RU      (152), VD      (152), RDU    (152), SMLR   (152)
8  DELR   ( 37), P      (152), P1      (152), PB      (152), PH1   (152)
  COMMON  P2      (152), SV      (152), RHO    (152), THETA (152)
1  W      (152), E      (152), EI      (152), EK      (152), A      (152)
2  V      (152), G      (152), D      (152), C      (152), X2     (152)
3  X3     (152), X4     (152), X5     (152), X6     (152), X7     (152)
4  SMLA   (152), SMLB   (152), SMLC   (152), SMLD   (152), SMLE   (152)
5  EC     (152), ER     (152), SMLQ   (152), SMLH   (152), RIGA   (152)
6  BIGB   (152), CV     (152), HC     (152), RR     (152), CHIC   (152)
7  CHIR   (152), CAPAC  (152), CAPAR  (152), CRTC   (152), CTRT   (152)
8  CRTPC  (152), GOFER  (152), FEW    (152), CAR    (152), OKLM   ( 37)
  COMMON  TELM    ( 37), EKLM    ( 37), ELM    ( 37), FCLM   ( 37)
1  FRLM   ( 37), WLM    ( 37), GLM    ( 37), AMASNO ( 37), CHRNO  ( 37)
2  ZP1    ( 37), ZP2    ( 37), SOLID  ( 37), ECHCK  ( 37), RK     (104)
3  RL     ( 37), RHOK   (104), RDK    (104), THETAK (104), TEMP   ( 16)
4  HEAD   ( 12), MAXL   , MAXLM

  *****
  COMMON /NAMEC/  RLF , RINC , RPUSH , RR      , DR
  COMMON /NAMEC/  KCNT , LCNT , MCNT , N      , INO , IL , IR
  COMMON /NAMEC/  VAL ( 10), RTABRO(150), RTABTH(150), RTABRD(150),
1  RHOTAB(150), THETAB(150), RDTAB (150)

  *****
  COMMON /SREZS/ IZSWT
  DOUBLE PRECISION R0,R1,R2,R3,R4,G1,G2,GS,VS,VOL,ZMS
  CALCULATE G(I),R(I),DELTAR(I),SV(I)
  IRM1=IR-1
  IRM2 = IR - 2
  ILP1=IL+1
  K=MIN0(KCNT,2)
80 IF(RTABRO(K).GT.R(IL).OR.K.EQ.KCNT) GO TO 90
  K=K+1
  GO TO 80
90 CONTINUE
  IF(VAL(6).EQ.2.) GO TO 120
  SPACE ZONING
  CALCULATE ZONE WIDTHS
  DELTAR(IL)=VAL(2)
  IF (ILP1.GE.IRM2) GO TO 930
  DO 92 I = ILP1,IRM2
  DELTAR(I)=DELTAR(I-1)*VAL(4)
92 CONTINUE
  DELTAR(IRM1) = VAL(3)
  DO 93 I = ILP1,IRM2

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      J = ILP1 + IRM2 - I
      WTR = FLOAT(IRM1 - J)/FLOAT(IRM1 - IL)
      DELTAR(J) = WTR*DELTAR(J) + (1.-WTR)*DELTAR(J+1)/VAL(4)
93  CONTINUE
C   CALCULATE ZONE BOUNDARIES
930  R1 = DBLE(R(IL))
      DO 94 I = ILP1,IRM1
      R1 = R1 + DBLE(DELTAR(I-1))
      R(I) = SNGL(R1)
94  CONTINUE
C   INITIALIZE LEFT BOUNDARY OF ZONING REGION
      I=IL
      R0=DBLE(R(IL))
104  R4=DBLE(RTAURO(K))
C   INITIALIZE ZONE I
105  GS = 0.00
      R1=R0
      R3=R0+DBLE(DELTAR(I))
C   SUBZONE CALCULATION
106  R2=R3
      IF(K.NE.KCNT) R2=DMIN1(R3,R4)
      GS=ZMS(R1,R2,K) + GS
      R1=R2
      IF(R2.LT.R4.OR.K.EQ.KCNT) GO TO 108
C   INITIALIZE NEW SECTION OF DENSITY PROFILE
      K=K+1
      R4=DBLE(RTABRO(K))
C   IS ZONE I COMPLETE
108  IF(R2.LT.R3) GO TO 106
C   ESTABLISH MASS, SPECIFIC VOLUME AND RIGHT BOUNDARY OF ZONE
      G(I)=SNGL(GS)
      VOL=R3-R0
      GO TO (114,111,112),IALPHA
111  VOL=VOL*(R3+R0)
      GO TO 114
112  VOL=VOL*(R3**2+R3*R0+R0**2)
114  SV(I)=SNGL(VOL/GS)
      I=I+1
      IF (I.EQ.IR) GO TO 131
C   GENERATE NEW ZONE
      R0=R3
      GO TO 105
C   MASS ZONING
120  CONTINUE
C   CALCULATE ZONE MASSES
      G(IL)=VAL(2)
      DO 121 I = ILP1,IRM2
      G(I)=G(I-1)*VAL(4)
121  CONTINUE
      G(IRM1) = VAL(3)
      DO 122 I = ILP1, IRM2
      J = ILP1 + IRM2 - I
      WTR = FLOAT(IRM1 - J)/FLOAT(IRM1 - IL)
      G(J) = WTR*G(J) + (1.-WTR)*G(J+1)/VAL(4)
122  CONTINUE
C   INITIALIZE LEFT BOUNDARY OF ZONING REGION
      R3=DBLE(R(IL))
      I=IL
      IF (K.LT.KCNT) GO TO 123
      G1 = 1.020
      GO TO 125
123  IF(R(IL).EQ.RTABRO(K-1).AND.IZSWT.EQ.(-1)) GO TO 124
      G1=ZMS(R3,DBLE(RTABRO(K)),K)
      GO TO 125
124  G1 = DBLE(EC(K-1))
C   INITIALIZE ZONE I
125  G2=DBLE(G(I))
      R2=R3

```

```

      VOL = 0.00
      SUBZONE CALCULATION
C 126 GS=DMIN1(G1,G2)
      R1=R2
      CALL ZRAD(R1,R2,GS,VS,K)
      G1=G1-GS
      G2=G2-GS
      VOL=VOL+VS
      IF (G1.GT.0.00) GO TO 127
      IF (K.LT.KCNT) GO TO 1260
      G1 = G2
      GO TO 127
C 1260 INITIALIZE NEW SECTION OF DENSITY PROFILE
      K = K + 1
      G1=DBLE(EC(K-1))
      IF (IZSWT.EQ.0) G1=ZMS(DBLE(RTAHKO(K-1)),DBLE(RTARRO(K)),K)
C 127 IS ZONE I COMPLETE
      IF (G2.GT.0.00) GO TO 126
C 127 ESTABLISH WIDTH, SPECIFIC VOLUME AND RIGHT BOUNDARY OF ZONE
      DELTAR(I)=SNGL(R2-R3)
      R3=R2
      SV(I)=SNGL(VOL)/G(I)
      I=I+1
      IF (I.EQ.IR) GO TO 130
      R(I)=SNGL(R3)
      GO TO 125
C 130 IS RIGHT REGION BOUNDARY PRESCRIBED
      IF (R(IR).EQ.1.E20) GO TO 150
      131 R2 = DBLE(R(IR))
      DO 135 I = ILP1,IRM1
      J = ILP1 + IRM1 - I
      WTR = FLOAT(IR-J)/FLOAT(IR-IL)
      R2 = R2 - DBLE(DELTAR(J))
      R(J) = WTR*R(J) + (1.-WTR)*SNGL(R2)
      135 CONTINUE
      GO TO 160
      150 R(IR) = SNGL(R3)
      160 RETURN
      END

```

```

      SUBROUTINE ZRAD(X1,X2,GG,VV,KK)
      COMMON LMDA(9405)
      EQUIVALENCE (S1,LMDA(136)),(IALPHA,LMDA(62))
C*****
      COMMON /NAMEC/      RLF      , RINC      , RPUSH , RH      DR
      COMMON /NAMEC/      KCNT      , LCNT      , MCNT  , N      INO      , IL , IR
      COMMON /NAMEC/ VAL      ( 10), RTABRO(150), RTABTH( 0), RTABRD(150),
1  RHOTAB(150), THETAB(150), RDTAB (150)
C
C      COMPUTE RIGHT BOUNDARY R2 AND VOLUME VS GIVEN LE      BOUNDARY R1
C      AND MASS GS OF A REGION IN WHICH DENSITY IS LINE    IN R**ALPHA
      COMMON /SREZS/IZSWT
      DOUBLE PRECISION R1,R2,GS,VS,C1,C2,X1,X2,GG,VV
      DOUBLE PRECISION C3, C4, C5, C6
      R1=X1
      GS=GG
      K=KK
      IF(K.GE.2) GO TO 5
      K=2
      GO TO 12
      5 IF (RTABRO(K).NE.RTABRO(K-1)) GO TO 10
      R2=R1
      VS = 0.00
      GO TO 60
      10 IF (IZSWT.EQ.(-1)) GO TO 12
      S = RHOTAB(K) - RHOTAB(K-1)
      IF(S.NE.0.)GO TO 30
      12 VS = GS/DBLE(RHOTAB(K-1))
      GO TO (15,20,25),IALPHA
      15 R2 = VS + R1
      GO TO 60
      20 R2 = DSQRT(VS + R1**2)
      GO TO 60
      25 R2 = (VS + R1**3)**(1.00/3.00)
      GO TO 60
      30 A1 = RTABRO(K-1)
      A2 = RTABRO(K)
      C1 = R1
      GO TO (34,32,33), IALPHA
      32 A1 = A1**2
      A2 = A2**2
      C1 = R1**2
      GO TO 34
      33 A1 = A1**3
      A2 = A2**3
      C1 = R1**3
      34 S = S/(A2 - A1)
      C3 = DBLE(A1 - RHOTAB(K-1)/S)
      C4 = GS*DBLE(2./S)
      C5 = C1 - C3
      C2 = C3
      C6 = C5**2 + C4
      IF (C6.GT.0.00) C2 = C3 + DSIGN(1.00,C5)*DSQRT(C6)
      GO TO (35,36,37), IALPHA
      35 R2 = C2
      GO TO 38
      36 R2 = DSQRT(C2)
      GO TO 38
      37 R2 = C2**(1.00/3.00)
      38 VS = C2 - C1
      60 X2=R2
      VV=VS
      RETURN
      END

```

```

C      DOUBLE PRECISION FUNCTION ZMS(R1,R2,KK)
C      COMPILED APRIL 15, 1966
C*****
COMMON /NAMEC/      RLF , RINC , RPUSH , RR , DR
COMMON /NAMEC/      KCNT , LCNT , MCNT , N , INO , IL, IR
COMMON /NAMEC/ VAL   ( 10), RTABRO(150), RTAITH(150), RTAHRD(150),
1  RHOTAB(150), THETAB(150), RDTAB (150)

C
C      FORMERLY NAMED SAV
C      COMPUTES MASS FOR A REGION WITH DENSITY LINLAR (IZSWT=0) OR
C      CONSTANT (IZSWT=-1) WITH RESPECT TO R**IALPHA
COMMON LMDA(9405)
EQUIVALENCE(S1,LMDA(136)),(IALPHA,LMDA(62)),(ALPHA,LMDA(79))
DOUBLE PRECISION R1,R2,ZMS,B1,B2,R1A,R2A
COMMON/SREZS/IZSWT
K=KK
IF(K.GE.2) GO TO 5
K=2
2 S=0.
GO TO 12
5 ZMS = 0.00
IF (RTABRO(K).EQ.RTABRO(K-1)) RETURN
IF (IZSWT.EQ.(-1)) GO TO 2
S = RHOTAB(K) - RHOTAB(K-1)
12 R1A = R1
R2A = R2
A1 = RTABRO(K-1)
A2 = RTABRO(K)
GO TO (25,15,20),IALPHA
15 R1A=R1A**2
R2A=R2A**2
A1 = A1**2
A2 = A2**2
GO TO 25
20 R1A=R1A**3
R2A=R2A**3
A1 = A1**3
A2 = A2**3
25 S = S/(A2-A1)
B1 = DBLE(RHOTAB(K-1) - S*A1)
B2 = DBLE(.5 * S)
ZMS = B1*(R2A-R1A) + B2*(R2A**2 - R1A**2)
IF(ZMS.GE.0.00) RETURN
S1=48.0025
CALL UNCLE
END

```



```

C *****
C COMMON /NAMEC/      KLF , RINC , RPUSH , RR      , DR
COMMON /NAMEC/      KCNT , LCNT , MCNT , N        , INO , IL , IP
COMMON /NAMEC/ VAL   ( 10), RTABRO(150), RTABTH(150), RTABRD(150),
1 RHOTAB(150), THETAB(150), RDTAB (150)
SOLVE QUADRATIC FOR COEFFICIENTS

E1 = A0 + A1*R1 + A2*R1**2
E2 = A0 + A1*R2 + A2*R2**2
EGS = KHO*A0*(R2**ALPHA-R1**ALPHA) +
      (ALPHA*RHO*A1)/(ALPHA+1.0)*(R2**(ALPHA+1.0)-R1**
      (ALPHA+1.0)) +
      (ALPHA*RHO*A2)/(ALPHA+2.0)*(R2**(ALPHA+2.0)-R1**(ALPHA+
      2.0))

ZS1(R2,R1) = R2-R1
ZS2(R2,R1) = R2 + R1
ZS3(R2,R1) = R2**2 + R1*R2 + R1**2
ZS4(R2,R1) = R2**3 + R1*R2**2 + R1**2*R2 + R1**3
ZS5(R2,R1) = R2**4+ R1*R2**3 + R1**2*R2**2 + R1**3*R2 + R1**4

C IRM1 = IR-1
EG = 0.0
K = IL

```



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```

DO 100 I = IL,IRM1
IF (E(I).NE.0.) GO TO 20
K = I + 1
GO TO 100
20 EG = EG + E(I)*G(I)
100 CONTINUE
IF (K.GE.IR-2) RETURN
E1 = E(K-1)
E2 = E(IR)
R1 = R(K)
R2 = R(IR)
105 CONTINUE
EGS = SV(K) * EG
C
C
C          CALCULATE COEFFICIENTS
C
B1 = ZS1(R2,R1)
GO TO (110,120,130),IALPHA
110 CONTINUE
C
B2 = (0.5*ZS2(R2,R1)-R2)
B3 = (ZS3(R2,R1)/3.0-R2**2)
A2 = ((EGS/B1-E2)-(E2-E1)*B2/B1)/(-ZS2(R2,R1)*B2+B3)
GO TO 140
C
120 CONTINUE
B2 = (2.0*ZS3(R2,R1)/(3.0*ZS2(R2,R1))-R2)
B3 = 0.5*ZS4(R2,R1)/ZS2(R2,R1)-R2**2
A2 = ((EGS/(B1*ZS2(R2,R1))-E2)-(E2-E1)*B2/B1)/(-ZS2(R2,R1)*B2+B3)
GO TO 140
C
130 CONTINUE
B2 = (0.75*ZS4(R2,R1)/ZS3(R2,R1)-R2)
B3 = (0.60*ZS5(R2,R1)/ZS4(R2,R1)-R2**2)
A2 = ((EGS/(B1*ZS3(R2,R1))-E2)-(E2-E1)*B2/B1)/(-ZS2(R2,R1)*B2+B3)
C
140 CONTINUE
A1 = (E2-E1)/B1-A2*ZS2(R2,R1)
A0 = E1-A1*R1-A2*R1**2
C
C
C          FIND THE ENERGY FOR EACH ZONE
C
DO 190 I = K, IRM1
R2 = R(I+1)
R1 = R(I)
C
GO TO (150,160,170),IALPHA
150 CONTINUE
C
E(I) = A0+0.5*A1*ZS2(R2,R1) + A2*ZS3(R2,R1)/3.0
GO TO 180
C
160 CONTINUE
E(I) = A0 + 2.0*A1*ZS3(R2,R1)/3.0 + A2*ZS4(R2,R1)/2.0
GO TO 180
C
170 CONTINUE
E(I) = A0 + 0.75*A1*ZS4(R2,R1) + 0.60*A2*ZS5(R2,R1)
180 CONTINUE
EI(I) = E(I)
ICU = I
IF (E(I).GE.E1) GO TO 190
GO TO 199
190 CONTINUE
RETURN

```

```
199 CONTINUE
    DO 200 J=K,ICU
        E(J)=E1
        EI(J)=E(J)
        EG=EG-E(J)*G(J)
200 CONTINUE
202 K = ICU + 1
    IF(ICU.EQ.IRM1) RETURN
    R1 = R(K)
    R2 = R(IR)
    GO TO 101
```

C  
END

## SECTION II

### FORTRAN ANALYSIS ROUTINE

#### 2.1. INTRODUCTION

The FORTRAN Analysis Routine is written entirely in FORTRAN IV for a 65K-UNIVAC 1108 computer. Conversion to other dialects of FORTRAN IV is not expected to present major obstacles. This routine consists of two independent programs. Program One is a variable renaming program; i.e., it provides for replacement of specified variable names by other names in a FORTRAN source deck, edits the deck by providing appropriate spacing, punches out a new sequenced deck, and provides a listing of this deck. In addition, it generates a tape containing all the required information for the other program. Program Two provides a printout of an extensive cross reference dictionary for the variables used in the program and the source line in which each variable appeared, including a description of the variable, its uses in each subprogram, and the subprogram in which it appears. It is intended that these programs be widely used for program maintenance.

#### 2.2. DESCRIPTION

In general, the input to Program One consists of an unlimited number of compilable FORTRAN IV subprograms, each of which is preceded by up to 15 "valid" strings which may appear in the subprograms that follow. For each "valid" string to be substituted for, a substitution string must be provided which maintains the compilability of the newly created statement.

A string is a sequence of characters. Only a valid string may be modified. To be valid, a string must be one of the following:

1. A variable name or a common block name or a subprogram name (hereafter the word "identifier" will be used to refer to one of these three),
2. A variable or subprogram name followed by a balanced set of parentheses,
3. The contents of a balanced set of parentheses,
4. Any string contained between consecutive commas,
5. Any string contained between a left parenthesis and the first comma that follows it, if any, prior to the corresponding right parenthesis,
6. Any string contained between the right-most comma following a left parenthesis, if any, and the corresponding right parenthesis;

and it must not be any of the following:

1. An arithmetic or logical operator,
2. A single non-alphabetic character,
3. .TRUE., .FALSE., .T., .F.,
4. A real or double precision or integer constant,
5. Any hollerith string (as appearing in a format statement or data statement or as an argument of a call).

In an arithmetic statement or arithmetic function statement, the left side of an equal sign is always a valid string; however, the right side may or may not be a valid string. The entire contents of the balanced set of parentheses following an IF statement is not necessarily a valid string.

In determining whether a substitution should take place in the subprograms, blanks are ignored. That is, all statements in the input programs are compressed as though no blanks were left anywhere. Of course, blanks are not deleted from hollerith fields.

### 2.3. EXAMPLES

1. A=B(1,I)+SQRT(D(2,2))-F(I,J+1,K)+2 \*G
2. IF(Z.LT..3) CALL EXIT

In the above two statements, only the following are valid strings:

* A	* F
E(1,I)	* I
* B	J+1
* I	* J
SQRT(D(2,2))	* K
* SQRT	* G
* D(2,2)	* Z
* D	* EXIT
F(I,J+1,K)	

Two passes are normally made through each statement in a program. A statement, including all its continuation cards, is read, and blanks are deleted. During the first pass every valid string is checked for substitution, and the substitution, if any, is made. During the second pass all identifiers and all variables having a completely numerical subscript are written on tape along with the new source line in which each variable appears. Those valid strings listed above with asterisks preceding them would have been recorded on tape. Not only are the name and source line of each identifier and numerically subscripted variable recorded on tape, but the fact that it appeared:

1. In a type statement or specification statement as (1) complex, (2) integer, (3) external, (4) logical, (5) real, (6) double precision, (7) equivalence, (8) parameter, (9) data.
2. On the left side of an equal sign.
3. As an argument of (1) the subprogram itself, (2) a referenced subroutine, (3) a referenced function.
4. As the name of the subprogram in a subroutine statement or function statement.

5. As the name of a referenced subroutine or referenced function.
6. In a statement which specifies a variable to be dimensioned (in a common or type statement with a numerical subscript or in a dimension statement), in which case its dimension and the number of characters in its dimension are also recorded.
7. In a common statement as either (1) a block common name or (2) a variable name, in which case its common block or blank is also recorded.
8. As a variable with a completely numerical subscript, in which case its subscript and the number of characters in its subscript are also recorded.
9. In a read statement.
10. Simply as a variable in none of the above.

This information is recorded at the time of appearance for each and every appearance of every identifier and numerically subscripted variable.

Pass two of Program One edits each statement after substitutions, if any are made, in the following manner: Single blanks are inserted before and after FORTRAN logical operators and the connectors =, +, and -, following a comma and neither before nor after the connectors \*, /, \*\*. Blanks do not occur in column 7 except for continuation cards, when they always occur. Identifier names are never split between one card and its continuation card. The continuation card mark in column 6 is always an apostrophe. The following statements are not edited except that the continuation marks are changed to apostrophes and information is placed in columns 73-80: comment cards, FORMAT statements, CONTINUE cards, and END statements. The compound or logical IF statement is split so that the second half of the compound statement begins on a continuation card. On all cards, columns 73-76 will contain the first four characters of the subprogram name appearing as the first name on the FOR card for each subprogram. Columns 77-80 will contain a 4 digit sequence number starting from 1 for each subprogram and increasing by 1.

A new sequenced source deck will be punched. A listing of the source deck will be printed with the additional information that the first line of each statement where a substitution has been made will be so noted. Any statement which required more than 19 continuation cards after editing will be so noted, and the additional cards will be punched and listed. All strings for which substitutions were requested that did not appear in the subprogram just processed will be noted at the end of the listing for that subprogram.

In addition to the pertinent description of Program One given above, the following unrelated facts concerning this program must be stated and understood if it is to be used properly:

1. Statement numbers are in no way affected by the program.
2. This program does not have extensive checking features, and a non-compilable statement may create an endless loop.
3. Each subprogram must begin with a FOR card with column one blank (the 7-8 punch removed) and end with an END statement.
4. Superfluous real, integer, and equivalence statements, created by virtue of variable substitutions, will not be eliminated.
5. No new real- or integer-type statements will be generated to correct for an unintentional change of type when a name substitution changes the type.
6. A name substitution will take place whenever that name appears regardless of whether it is subscripted.
7. A variable plus subscript substitution will take place only when the variable and subscript agree exactly. However, no substitution will take place if the subscript is numeric and appears in a dimension statement or any statement which specifies its dimensionality. Thus, the replacement of R(152) by PCYCLE would not occur in the statement COMMON R(152).
8. The appearance of any variable followed by a left parenthesis not appearing in a specification statement defining its dimensionality will be checked to determine whether it has been



previously dimensioned. If not, it will be assumed to be a function reference. Thus, on substituting C(5) for B, if C has not been previously dimensioned, it will be considered a function and will not be otherwise noted as a possible source error.

9. This program cannot accept comment cards embedded between the first card of a statement and the last continuation card of the statement. Such cards must be removed from a deck before this program is used. Comment cards may appear any place else between a FOR card and an END card of a subprogram.
10. Anything appearing in columns 73-80 on comment cards is deleted, and the sequencing scheme previously described is substituted.
11. The number of + signs, - signs, and logical operators appearing in a compound or logical IF statement must not exceed 20 if this statement is to be reproduced correctly.

Program Two processes the output of Program One. It will print out the following information for each subprogram:

1. The subprogram name plus an 18 character tape identification. This identification is described further in Section 2.4 of this report.
2. An alphabetic listing of all common blocks starting with blank common. For each common block the following information is printed.
  - a. A listing of the variables within each block in order of their appearance in the block, and for each variable its octal and decimal origin relative to zero within the block.
  - b. An alphabetic listing of the variables within each block, and for each variable its octal and decimal origin relative to zero within the block.



3. An alphabetic listing of the subroutines and functions referenced within this subprogram and the source lines in which they were referenced.
4. An alphabetic listing of each variable appearing in the subprogram. For each variable the following information is printed:
  - a. A single line containing the variable name, the block name in which it may appear, its octal and decimal origin relative to zero within that block, its dimensions (if any), the size of its dimensions, and a series of code letters describing its "uses" within the program (i. e. , its type, whether it appears anywhere within the subprogram on the left side of an equal sign, whether it appears anywhere within the subprogram as the argument of a call, whether it appears anywhere in the subprogram in a read statement, etc.) A full description and listing of these code letters appears in Section 2.4 of this report and on the first page of the printed output of Program Two.
  - b. An ordered listing of every source line in which the variable appears is printed. Following each source line there may appear the same coded letters as are described in (4a), above, if the variable can be modified within that line (by appearing on the left side of an equal sign or as an argument of a function, in a read statement, etc.).
  - c. For each variable that is numerically subscripted within the program, a sort order listing for each numeral subscript, followed by a list of all source lines in which it appears with this subscript.

After all subprograms have been processed and each of the above listings has been printed, the program prints an alphabetic listing of all

3. An alphabetic listing of the subroutines and functions referenced within this subprogram and the source lines in which they were referenced.
4. An alphabetic listing of each variable appearing in the subprogram. For each variable the following information is printed:
  - a. A single line containing the variable name, the block name in which it may appear, its octal and decimal origin relative to zero within that block, its dimensions (if any), the size of its dimensions, and a series of code letters describing its "uses" within the program (i. e., its type, whether it appears anywhere within the subprogram on the left side of an equal sign, whether it appears anywhere within the subprogram as the argument of a call, whether it appears anywhere in the subprogram in a read statement, etc.) A full description and listing of these code letters appears in Section 2.4 of this report and on the first page of the printed output of Program Two.
  - b. An ordered listing of every source line in which the variable appears is printed. Following each source line there may appear the same coded letters as are described in (4a), above, if the variable can be modified within that line (by appearing on the left side of an equal sign or as an argument of a function, in a read statement, etc.).
  - c. For each variable that is numerically subscripted within the program, a sort order listing for each numeral subscript, followed by a list of all source lines in which it appears with this subscript.

After all subprograms have been processed and each of the above listings has been printed, the program prints an alphabetic listing of all

identifiers appearing in any of the subprograms, and for each identifier, an alphabetic listing of the subprograms in which it appeared. The subprograms are listed by subroutine name (not by the name on the FOR card). This ends the output description of Program Two.

Both Program One and Program Two use logical unit 10 to store and process the identifier information. A complete description of the format and information stored on logical unit 10 by Program One is contained in Section 2.5 of this report. Table lengths for Program Two are also given in that section.

#### 2.4. INPUT

It is suggested that FORTRAN code sheets be used for all data input, since the data format has been so designed. There are three types of data input for Program One. Data type 1 is a single card having two fields. It precedes a complete run of Program One which may consist of one or more subprograms. Field 1 - Column 1 (format 11) controls logical unit 10; if field 1 is blank, logical unit 10 will be rewound. This implies either that the unit is blank or that information previously stored on this unit is no longer desired. If field 1 contains a one, it is assumed that logical unit 10 had information previously stored on it from a previous run of Program One; then logical unit 10 will be positioned forward to the end of the previously stored information, ready to receive additional information. Field 1 - Columns 2-6 inclusive are blank. Field 2 - Columns 7-74 (format 3A6) is the secondary identification valid for all subprograms of this run. It may contain any information desired, e.g., the date. It is intended that a program will be written to edit logical unit 10, deleting one version of a subprogram and accepting another version of this same subprogram by utilizing the uniqueness of the 18 character identification in field 2 between runs of Program One.

For Data type 2, one or more cards must precede each subprogram, but no subprogram will require more than 15 of these cards. Fifteen "blocks" have been set aside, internal to Program One, for "substitutions." (A "substitution" is defined as the combination of a source program string to be changed and its replacement string.) A block in which no substitution has ever been placed or one in which a substitution has been deleted without replacement is considered void. A block that is not void is full. To "delete" means to wipe out a substitution. It can only be brought back by fully restating the substitution on a new type 2 card, as described below. A type 2 card may serve any of the following purposes:

1. It may fill a void block with a substitution. The block is then full. The substitution then applies through all subprograms that follow it unless deleted or replaced by another type 2 card.
2. It may delete a particular full block. That block is then void.
3. It may delete all existing full blocks. All blocks will then be void.
4. It may delete a substitution from a full block and replace it with another substitution in the same block. That block is then still full. The new substitution will then apply through all subprograms that follow it unless it is specifically deleted or replaced by still another type 2 card.

Type 2 cards consist of three fields, hereafter referred to as F1, F2, and F3. F1 (Format A6) delineates the last data type 2 card by having some non-blank character in any of columns 1-6 of the data card, (except that column 1 must not have a 7-8 punch). That is, columns 1-6 must be blank for all but the last data type 2 card preceding each subprogram.

F2 (format 71A1) is either entirely blank, in which case it sets the block specified in F3 void, or contains the substitution in columns 7-77, inclusive. If non-blank, the general form of F2 is SOURCE PROGRAM STRING TO BE MODIFIED = ITS REPLACEMENT STRING. The following rules must be observed in the specification of F2.

1. F2 must start in column 7. (It is considered blank if column 7 of F2 is blank.)
2. The two strings must be separated by an equal sign.
3. There must not be any blanks between column 7 and the last character of the replacement string.
4. Column 78 must be blank.

For F3 (Format I2), columns 79 and 80 must contain blanks or the numbers 01-15 or the number 99. F3 may be blank if and only if F1 is non-blank and F2 is blank, i.e., to express the end of type 2 data without supplying any information on F2 and F3. In general, F3 contains the block numbers 01-15 for which an F2 substitution or deletion or replacement substitution is being made. The information for a block need not be ordered, i.e., information for block 07 may be supplied before information for block 01, and blocks may be arbitrarily full or void. If F3 contains a 99, all 15 blocks are void. F2 would be blank if F3 contained a 99, and F1 might or might not be blank. It is perfectly valid for additional type 2 cards to follow an F3 containing a 99 (in which case F1 would be blank when F3 contained a 99).

If all 15 blocks are void upon return from the program that processes the type 2 cards, then Program One will bypass its first pass (the substitution phase) and will presumably run twice as fast. This implies that all one is interested in is the cross reference dictionary generated in pass two. If this is so, then the data type 2 cards preceding each subprogram should consist of a single card with perhaps a single character in column 1.

(However, the single character must not be a 7-8 punch, which transfers control to the 1108 operating system.)

Data type 3 cards are subprogram source cards to be processed. They must start with a FOR card with the column 1 blank (that is the 7-8 punch must be removed). Following column 1 there must be at least one blank preceding and at least one blank following the "FOR" and before its Deck Name. Just prior to the FOR card there must be a data type 2 card with at least one character other than a blank in columns 1-6 (and not a 7-8 punch in column 1). Following the modified FOR card are the source cards, which must be compilable. The source deck must end with an END card.

Program Two requires no card input. A request for logical unit 10 (ASG) must be supplied by the user for both Program One and Program Two.

## 2.5. APPENDIX: PROGRAM DETAILS

Table I contains information describing the code letters which define the "use" of an identifier (see steps (4a) and (4b) of the Program Two output description in Section 2.2), and the code numbers written on tape by Program One which identify that "use."

Any combination of the code letters appearing in column 1 of Table I may be printed in the Program Two output previously described in step (4a) of Section 2.3, but only those code letters in Table I which have an asterisk preceding them will be printed out as described in step (4b). The latter code letters reference only those statements which can modify a variable.

The information written by Program One on logical unit 10 is written in unformatted I/O (equivalent to binary) and consists of six words per record. The six words make up four fields. Field 1 is in field data and is either (1) the program name appearing on the FOR card, as the first word of the first and last record of each subprogram, or (2) an identifier appear-

Table I

## DESCRIPTION OF CODE LETTERS

<u>Code Letter</u>	<u>Code Number(k)</u>	<u>Description</u>
U	- 89	Variable appears in a <u>parameter</u> statement
* D	- 90	Variable appears in a <u>data</u> statement
	- 91	Name appears on <u>FOR</u> Card
Z	- 92	Variable appears in a <u>complex</u> type statement
* X	- 93	Variable appears on <u>left side of an equal sign</u>
* H	- 94	<u>Variable</u> appears in a <u>common</u> block
	- 95	<u>Block name</u> appears in a <u>common</u> statement
* S	- 96	Variable appears as an <u>argument of a call</u>
B	- 97	Variable appears in a <u>double precision</u> type statement
* R	- 98	Variable appears in a <u>read</u> statement
	- 99	<u>Subroutine name</u> on subroutine or <u>function statement card</u>
E	-100	Variable appears in an <u>external</u> type statement
C	-101	<u>Subroutine name</u> appears in a <u>call</u> statement
i	-102	Variable name appears in an <u>integer</u> type statement
* Q	-103	Variable name appears in an <u>equivalence</u> type statement
L	-104	Variable appears in a <u>logical</u> type statement
F	-105	<u>Function name</u> appear in an arithmetic statement or <u>if</u> statement, or on either side of the equal sign in an arithmetic function statement (but not on a function statement card described in -99 above)
P	-106	Variable appears in a <u>real</u> type statement
A	-107	Variable appears as a <u>program argument</u> on a function or subroutine statement card
* G	-108	Variable appears as the <u>argument of a referenced function</u> as described in -105 above
	-18 ≤ k < 0	Variable appears in a <u>dimension</u> statement or in a statement describing its dimensionality where  k  is the number of characters in its dimension description including parentheses, commas, and digits
	0 < k ≤ 18	Variable appears in a statement, other than one describing its dimensionality or an equivalence statement, with a <u>numerical subscript</u> where k is the number of characters in its subscript including both parentheses, commas, and digits
	-218 ≤ k < -200	Variable appears in an <u>equivalence</u> statement with a subscript where  k-200  is the number of characters in its subscript, including both parentheses, commas, and digits
	0	Not specified above (as the variables A or B in the arithmetic statement Y=A+B)



ing in the program. Field 1 is left adjusted and filled with blanks on the right if the identifier is less than six characters.

Field 2 is in field data and consists of three words, either:

1. The 18 character identification specified on the data type 1 card of a specific run of Program One. It will appear in the second field of the first record of every subprogram processed during this run.
2. A variable dimension, left adjusted, with no embedded blanks, containing both parentheses and their contents, and filled with blanks on the right. Field 2 will contain this information only if the variable in field 1 is appearing in a statement specifying its dimensionality at the time this record is written.
3. A variable numerical subscript, left adjusted, with no embedded blanks (as described in step 2 above). Field 2 will contain this information only if the variable in field 1 is appearing in any statement other than one specifying its dimensionality at the time this record is written and is numerically subscripted.
4. A block name (which may be blank for blank common). Word 1 of field 2 will contain this information only if the variable is appearing in a common statement at the time the record is written; then words 2 and 3 of field 2 will be blank. If field 1 contains the block name, the first word of field 2 will also contain that block name. All three words of field 2 contain blanks.

Field 3, the fifth word of the record, is the code number appearing in the second column of Table I. The last record of each subprogram contains a -200 in field 3 (generated by the END statement). Field 4 contains the source line number in which the identifier in field 1 appears. Following the last record of the last subprogram is one additional six word record containing a -200 in its fifth word. There is no end-of-file following this.



One can, therefore, determine the end of information by seeking two consecutive records each of which contains -200 in its fifth word.

If a variable appears in a statement specifying its dimensionality or is numerically subscripted, it will appear in two consecutive records. For example, the numerically subscripted variable A(10) in the statement, A(10)=Y appearing in line 17 of a subprogram appears in two records as:

	<u>Field 1</u>	<u>Field 2</u>	<u>Field 3</u>	<u>Field 4</u>
Record n	A	(10)	4	17
Record n+1	A		-93	17
Record n+2	Y		0	17

The purpose of the above description is to provide the user with an opportunity to write his own specialized cross reference dictionary (i.e., his own variation of Program Two. For example, he may wish to print out all variables appearing in READ statements throughout the program in the order in which they are read.

Certain assumptions had to be made in Program Two concerning the internal allocation of space for the many tables required by the cross reference dictionary. This allocation is further complicated by a possible variation in core space that will be available to the user in the near future. It is currently assumed that the total data space must not exceed 40,000. That is, the sum of the "total space" created by the parameter occurrences listed in Table II and fixed working memory must not exceed 40,000. However, as long as the total does not exceed 40,000, any one or more of the five parameters may be modified within the program in the event the individual tables to which they refer are exceeded. This requires the change of a single PARAMETER card in Program Two, which states: PARAMETER Z1=15000, Z2=500, Z3=1000, Z4=2500, Z5=1000, where Z1-Z5 and the number of appearances of Z1-Z5 are as defined in Table II. The "limit" column in Table II refers to the maximum appearances or references or number of such items in any one subprogram.

Table II  
DEFINITION OF ENTRIES IN PARAMETER CARD

<u>Para- meter</u>	<u>Description</u>	<u>Limit</u>	Number of tables requiring this para- meter =			<u>Total space</u>
Z1	Number of references of variables with or without subscripts	15000	×	1	=	15,000
Z2	Number of unique common block names	500	×	1	=	500
Z3	Number of appearances of variables with unique subscripts	1000	×	6	=	6,000
Z4	Number of appearances of unique identifiers	2500	×	4	=	10,000
Z5	Number of appearances of unique variables that are dimensioned	1000	×	5	=	5,000
						36,500
	Fixed working memory (not subject to change)					3,500
						40,000

SECTION III  
DATA-LINK EDITING ROUTINES

3.1. INTRODUCTION

In order to facilitate use of the computer codes developed in the NEIS on the CDC-6600 computer at the Air Force Weapons Laboratory, a data link was installed between the SNEL facility in San Diego and the AFWL facilities in Albuquerque. The link terminals include an IBM-1978 reader-printer at SNEL, which communicates with an IBM-7702 tape terminal at AFWL. The format of the information which can be transmitted or received by the 1978 is unfortunately incompatible with the requirements of the Chippewa operating system for the 6600 computer, so editing procedures are required as an intermediate step in the data transmission process. Two programs, called TOGA and FROGA, were written for the AFWL CDC-160A computer to perform these editing functions.

3.2. TOGA

The TOGA program edits BCD output tapes prepared by the 6600, producing modified tapes which can be read by the 7702, and printed on the 1978. The following editing functions are accomplished:

1. Printer carriage control characters written in FORTRAN form are blank, 0, -, and 1, indicating single space, double space, triple space, and page eject. These actions precede the printing of the line in which they occur. The corresponding 1978 control characters are /, S, T and A, respectively, with the difference that the actions follow the printing of the line in which they occur. The TOGA code therefore replaces each FORTRAN carriage control character by its 1978 counterpart, which is then placed

at the head of the preceding line. If the first character in the line is already a 1978 character, no action is taken; if it is an illegal character, a single space (/) is supplied.

2. The 1978 requires each line to be terminated by a segment mark, which is a special BCD character. This character is placed at the end of each line, after any trailing blanks are deleted.
3. The physical records transmitted over the link are blocked to some extent. The number of lines per block varies from two to seven, depending upon the total number of characters, which may not exceed 328.
4. The length of a line printable on the 1978 is 120 characters, plus the leading control character and the trailing segment mark. If the input lines exceed this length, the overflow goes onto a second line flagged by a leading exclamation point, following which the carriage is double-spaced.
5. Input records less than three characters long are considered noise and are skipped.

### 3.3. FROGA

The FROGA program edits information transmitted by the 1978 into a form acceptable for job input by the Chippewa system. The following functions are performed:

1. The transmitted records, except possibly the last, consist of four 80 character card images, each followed by a segment mark. These records are read one at a time. For each card image, the segment mark and any trailing blanks are deleted, the latter two at a time.
2. The card images are moved to an output area. The end of each card image is indicated by the following rule: The number of

characters for each card image must be a multiple of 10, and the last two of these must be 6-bit (binary) zeros.

3. As the BCD characters are moved to the output area, they are translated into the Chippewa display code by table lookup.
4. The size of the output buffer is 5120 characters. The buffer is written onto an output "display tape" when full, or when an end-of-record flag is received. The end-of-record flag is an apostrophe in column 1 of a card image. If the buffer is just filled before the end-of-record flag is received (an unlikely occurrence), a special short record consisting of eight binary zeros (48 bits) is written. Either this special short record or the normal short record (less than 512 ten character words) indicates to the Chippewa system that a "logical record" is complete.
5. End-of-file records are written on the display tape when a second apostrophe is received following the first, e.g., in column 2 of the same card image. This end-of-file separates one job from another. A third apostrophe indicates end of transmission, as well as end-of-record and end-of-file. When only one job is transmitted at a time, only single or triple apostrophe flags are used (the preferred procedure). On end-of-transmission, two end-of-file records are written on the display tape.

characters for each card image must be a multiple of 10, and the last two of these must be 6-bit (binary) zeros.

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